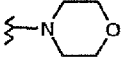
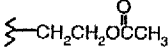
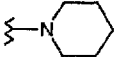


Example #	R <sub>1</sub>	Name	NMR Data*	Mass Spec.
48		2,2'-[[4-(4-Morpholinyl)phenyl]methylene]bis[4-[[[(5-methyl-1H-tetrazol-1-yl)imino]methyl]]phenol	<sup>1</sup> H NMR in DMSO: 10.47(s, 2H); 9.15(s, 2H); 7.78(dd, J=1.76, 2.34, 8.50Hz, 2H); 7.46(d, J=1.76Hz, 2H); 6.94(m, 6H); 5.99(s, 1H); 3.73(t, J=4.69Hz, 4H); 3.07(t, J=4.69Hz, 4H); 2.47(s, 6H)	(M+H) <sup>+</sup> 580.1
49		4-[[Bis[[[5-(5-methyl-1H-tetrazol-1-yl)imino]methyl]-2-hydroxyphenyl]methylene]benzeneethanol, acetate ester	<sup>1</sup> H NMR in DMSO: 5.20(s, 2H); 9.15(s, 2H); 7.80(dd, J = 1.7, 8.2, 8.8Hz, 2H); 7.45(d, J = 1.7Hz, 2H); 7.03(d, J = 8.2Hz, 2H); 7.03(d, J = 8.2Hz, 2H); 6.05(s, 1H); 4.20(t, J = 6.4, 7.0Hz, 2H); 2.46(s, 3H); 1.98(s, 3H)	580, 552, 525, 456
50		2,2'-[[4-(1-Piperidinyl)phenyl]methylene]bis[4-[[[(5-methyl-1H-tetrazol-1-yl)imino]methyl]]phenol	<sup>1</sup> H NMR in DMSO: 10.48(s, 2H); 9.15(s, 2H); 7.80(d, J=8.21 Hz, 2H); 7.47(s, 2H); 7.00(d, J=8.79 Hz, 2H); 6.93(s, 2H); 5.98(s, 1H); 3.11(m, 4H); 2.47(s, 6H); 1.62(m, 4H); 1.54(m, 2H)	576 (M-H)

\*All <sup>1</sup>H NMR and <sup>13</sup>C NMR spectra were acquired on a Varian Mercury VX 300 Spectrometer and referenced to tetramethylsilane (TMS) unless indicated otherwise. Chemical shifts and coupling constants are reported in parts per million (ppm) and Hertz (Hz) respectively.